

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN073020\
 Data File : BN011299.D
 Acq On : 30 Jul 2020 14:53
 Operator : CG/JU
 Sample : SSTD1.604
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.604

Quant Time: Jul 30 15:25:10 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 30 14:13:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	1229	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	4251	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2666	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	6027	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	5561	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	4918	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	17270	1.66	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	46235	1.78	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.26	128	19742	1.552	ng/ul	99
5) 2-Methylnaphthalene	11.89	142	13663	1.643	ng/ul	98
7) Acenaphthylene	13.82	152	20607	1.595	ng/ul	97
8) Acenaphthene	14.16	153	16002	1.576	ng/ul	99
9) Fluorene	15.16	166	19074	1.655	ng/ul	98
11) Pentachlorophenol	16.52	266	2444	1.898	ng/ul	99
12) Phenanthrene	16.90	178	31894	1.606	ng/ul	99
13) Anthracene	16.99	178	29108	1.778	ng/ul	99
15) Fluoranthene	18.93	202	39255	1.769	ng/ul	100
17) Pyrene	19.29	202	38315	1.290	ng/ul	97
18) Benzo(a)anthracene	21.06	228	34825	1.771	ng/ul	99
19) Chrysene	21.11	228	37412	1.431	ng/ul	100
21) Benzo(b)fluoranthene	22.57	252	33571	1.581	ng/ul	94
22) Benzo(k)fluoranthene	22.62	252	36099	1.481	ng/ul#	90
23) Benzo(a)pyrene	23.11	252	28615	1.602	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.27	276	30655	1.456	ng/ul	100
25) Dibenzo(a,h)anthracene	25.29	278	24427	1.538	ng/ul	94
26) Benzo(g,h,i)perylene	25.90	276	26862	1.284	ng/ul	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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